

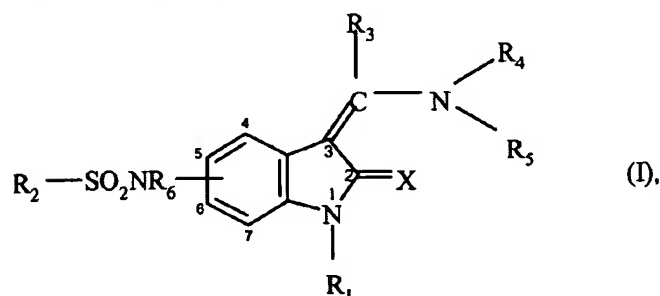
CLAIMS

Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

Claim 1 (currently amended): A compound of the formula I



or a pharmaceutically acceptable salt thereof, wherein:

X is an oxygen or sulphur atom,

R₁ is a hydrogen atom, a C₁₋₄-alkoxycarbonyl or C₂₋₄-alkanoyl group,

R₂ is a C₁₋₆-alkyl group optionally substituted by one or more halogen atoms or a phenyl group or a C₂₋₆-alkenyl group optionally substituted by a phenyl group, wherein the phenyl moiety may be substituted in each case by a fluorine, chlorine, bromine or iodine atom, by a C₁₋₃-alkyl or C₁₋₃-alkoxy group,

a phenyl group which may be mono- or disubstituted by fluorine, chlorine, bromine or iodine atoms, by C₁₋₃-alkyl or C₁₋₃-alkoxy groups, wherein the substituents may be identical or different,

a phenyl group substituted by a trifluoromethyl, carboxy, C₁₋₃-alkoxycarbonyl, aminocarbonyl, cyano, aminomethyl, nitro or amino group,

a C₄₋₆-alkyl, C₃₋₇-cycloalkyl, trimethylphenyl or naphthyl group,

~~a 5-membered heteroaromatic group optionally substituted by a C₁₋₃-alkyl group, which contains, in the heteroaromatic moiety,~~
~~an imino group optionally substituted by a C₁₋₃-alkyl group, an oxygen or sulphur atom,~~
~~an imino group optionally substituted by a C₁₋₃-alkyl group and an oxygen, sulphur or nitrogen atom,~~
~~an imino group optionally substituted by a C₁₋₃-alkyl group and two nitrogen atoms, or~~
~~an oxygen or sulphur atom and two nitrogen atoms, and to which a phenyl ring may be fused via two adjacent carbon atoms,~~
~~or is a 6-membered heteroaromatic group optionally substituted by a C₁₋₃-alkyl group, which contains one or two heteroatoms in the heteroaromatic moiety and to which a phenyl ring may be fused via two adjacent carbon atoms,~~

R₃ is a hydrogen atom or a C₁₋₆-alkyl group,

a phenyl group optionally substituted by a fluorine, chlorine or bromine atom, by a C₁₋₃-alkyl, hydroxy, C₁₋₃-alkoxy, C₁₋₃-alkylsulphenyl, C₁₋₃-alkylsulphinyl, C₁₋₃-alkylsulphonyl, phenylsulphenyl, phenylsulphinyl, phenylsulphonyl, nitro, amino, C₁₋₃-alkylamino, di-(C₁₋₃-alkyl)-amino, C₂₋₅-alkanoylamino or N-(C₁₋₃-alkylamino)-C₂₋₅-alkanoylamino group,

R₄ is a phenyl or naphthyl group ~~optionally~~ substituted by R₇, which may additionally be substituted by a chlorine or bromine atom or a nitro group, ~~a 5-membered heteroaromatic group which contains an imino group, an oxygen or sulphur atom or an imino group, an oxygen or sulphur atom and one or two nitrogen atoms, or~~
~~a 6-membered heteroaromatic group which contains one, two or three nitrogen atoms, while the abovementioned 5- and 6-membered heteroaromatic groups may additionally be substituted by a chlorine or bromine atom or by a methyl group or wherein a phenyl ring may be fused to the abovementioned 5- and 6-membered heteroaromatic groups via 2 adjacent carbon atoms, or~~

R₅ and R₆ in each case independently of one another are hydrogen atoms or C₁₋₃-alkyl groups, and

R₇ is a ~~fluorine, chlorine, bromine or iodine atom or a cyano group,~~
~~a methoxy group or a C₂₋₃-alkoxy group, which may be substituted in the 2 or 3 position by~~
~~an amino, C₁₋₃-alkylamino, di-(C₁₋₃-alkyl)-amino or 5- to 7-membered cycloalkyleneimino~~

group, while in each case an alkyl moiety in the abovementioned alkylamino and dialkylamino groups may additionally be substituted by a phenyl group, a trifluoromethyl, nitro, amino, C₁₋₃-alkylamino, di (C₁₋₃-alkyl)-amino, C₂₋₅-alkanoylamino, N-(C₁₋₃-alkyl)-C₂₋₅-alkanoylamino, C₁₋₅-alkylsulphonylamino, N-(C₁₋₃-alkyl)-C₁₋₅-alkylsulphonylamino, phenylsulphonylamino, N-(C₁₋₃-alkyl)-phenylsulphonylamino, aminosulphonyl, C₁₋₃-alkylaminosulphonyl or di (C₁₋₃-alkyl)-aminosulphonyl group, while in each case an alkyl moiety in the abovementioned alkylamino and dialkylamino groups may additionally be substituted by a carboxy, C₁₋₃-alkoxycarbonyl, aminocarbonyl, C₁₋₃-alkylaminocarbonyl, di (C₁₋₃-alkyl)-aminocarbonyl, 2-dimethylaminoethylaminocarbonyl or N-methyl-(2-dimethylaminoethyl)-aminocarbonyl group and in each case the alkyl moiety of the abovementioned alkanoylamino or alkylsulphonylamino groups may additionally be substituted by a phenyl, amino, C₁₋₃-alkylamino, di (C₁₋₃-alkyl)-amino or a 4- to 7-membered cycloalkylencimino group, a C₂₋₄-alkylamino group which is terminally substituted in the 2, 3- or 4 position by an amino, C₁₋₃-alkylamino, di (C₁₋₃-alkyl)-amino, benzylamino, N-(C₁₋₃-alkyl)-benzylamino, C₂₋₅-alkanoylamino or N-(C₁₋₃-alkyl)-C₂₋₅-alkanoylamino group and wherein additionally the amino-hydrogen atom may be replaced by a C₂₋₅-alkanoyl, benzoyl, C₁₋₅-alkylsulphonyl or phenylsulphonyl group, while the last-mentioned C₂₋₅-alkanoyl or C₁₋₅-alkylsulphonyl groups in the alkyl moiety may be substituted by a phenyl group, a carbonyl group which is substituted by a hydroxy, C₁₋₃-alkoxy, amino, C₁₋₃-alkylamino, N-(C₁₋₅-alkyl)-C₁₋₃-alkylamino or C₅₋₇-cycloalkylencimino group, a C₁₋₃-alkyl group which may be substituted by an amino, C₁₋₅-alkylamino, C₅₋₇-cycloalkylamino or phenyl-C₁₋₃-alkylamino group which may additionally be substituted at the amino-nitrogen atom in each case by a C₁₋₄-alkyl, C₅₋₇-cycloalkyl or C₂₋₄-alkenyl or C₁₋₄-alkyl group, while the abovementioned C₁₋₄-alkyl substituent in each case may additionally be mono-, di- or trisubstituted by a cyano, carboxy, C₁₋₃-alkoxycarbonyl, C₂₋₄-alkanoyl, pyridyl, imidazolyl, benzo[1,3]dioxol or phenyl group, while the phenyl group may be substituted by fluorine, chlorine or bromine atoms, by methyl, methoxy, trifluoromethyl, cyano or nitro groups and the substituents may be identical or different, or in the 2, 3- or 4 position by a hydroxy group,

a C₁₋₃-alkyl group which is substituted by a ~~hydroxy, carboxy, morpholino, thiomorpholino, 1-oxo-thiomorpholino, 1,1-dioxo-thiomorpholino, piperazine, N-(C₁₋₃-alkyl)-piperazine or N-benzyl-piperazine group~~, by a 5- to 7-membered ~~cycloalkenyleneimino group~~ or by a 4- to 7-membered ~~cycloalkyleneimino~~ piperidino group, while the abovementioned 5- to 7-membered ~~cycloalkyleneimino groups~~ piperidino group may be substituted by one or two C₁₋₃-alkyl groups, which may in turn be terminally substituted by a hydroxy, amino or C₂₋₄-alkanoylamino group, or by a C₅₋₇-cycloalkyl or phenyl group and by a hydroxy group and in the abovementioned ~~cycloalkyleneimino groups~~ piperidino group a methylene group adjacent to the nitrogen atom may be replaced by a carbonyl group;
~~a C₁₋₃-alkyl group which is substituted by a 5- to 7-membered cycloalkyleneimino group,~~
~~while a phenyl group optionally mono- or disubstituted by fluorine, chlorine or bromine atoms or by methyl or methoxy groups, wherein the substituents may be identical or different,~~
~~or an oxazolo, imidazolo, thiazolo, pyridino, pyrazino or pyrimidino group optionally substituted by a fluorine, chlorine, bromine or iodine atom, by a methyl, methoxy or amino group is fused to the abovementioned 5- to 7-membered cycloalkyleneimino groups via 2 adjacent carbon atoms, while the abovementioned monosubstituted phenyl groups may additionally be substituted by a fluorine, chlorine or bromine atom, by a methyl, methoxy or nitro group, or~~
is an imidazolyl or 1H-C₁₋₃-alkylimidazolyl group.

Claim 2 (original): A compound of formula I according to claim 1 wherein the sulphonylamino group of the formula R₂-SO₂NR₆- is linked to the 5-position of the indolinone group.

Claim 3 (original): A compound of formula I according to claim 1, wherein:

R₃ is a phenyl group optionally substituted by a fluorine, chlorine or bromine atom, by a C₁₋₃-alkyl, hydroxy, C₁₋₃-alkoxy, C₁₋₃-alkylsulphenyl, C₁₋₃-alkylsulphinyl, C₁₋₃-alkylsulphonyl,

phenylsulphenyl, phenylsulphanyl, phenylsulphonyl, nitro, amino, C₁₋₃-alkylamino, di-(C₁₋₃-alkyl)-amino, C₂₋₅-alkanoylamino or N-(C₁₋₃-alkylamino)-C₂₋₅-alkanoylamino group.

Claim 4 (original): A compound of formula I according to claim 1, wherein:

R₂ is a C₁₋₃-alkyl group optionally substituted by one or more halogen atoms or a phenyl group or a C₂₋₄-alkenyl group optionally substituted by a phenyl group, wherein the phenyl moiety in each case may be substituted by a fluorine, chlorine, bromine or iodine atom or by a C₁₋₃-alkyl or C₁₋₃-alkoxy group.

Claim 5 (currently amended): A compound of formula I according to claim 1, wherein:

X is an oxygen atom,

R₁ is a hydrogen atom,

R₂ is a C₁₋₃-alkyl group optionally substituted by one or more fluorine atoms or a phenyl group or a C₂₋₄-alkenyl group optionally substituted by a phenyl group;
a phenyl group which may be mono- or disubstituted by fluorine, chlorine, bromine or iodine atoms, by C₁₋₃-alkyl or C₁₋₃-alkoxy groups, wherein the substituents may be identical or different,

a phenyl group substituted by a trifluoromethyl, carboxy, C₁₋₃-alkoxycarbonyl, aminocarbonyl, cyano, aminomethyl, nitro or amino group, or

a C₄₋₆-alkyl, C₃₋₇-cycloalkyl, trimethylphenyl or naphthyl group, ~~or~~

~~a pyridinyl, quinolyl, isoquinolyl, oxazolyl, isoxazolyl, imidazolyl or 1-(C₁₋₃-alkyl)-imidazolyl group optionally substituted by a C₁₋₃-alkyl group;~~

R₃ is a hydrogen atom or a C₁₋₄-alkyl group, or

a phenyl group optionally substituted by a fluorine, chlorine, bromine or iodine atom, by a C₁₋₃-alkyl, C₁₋₃-alkoxy, nitro or amino group,

R₄ is a phenyl group ~~optionally~~ substituted by R₇,

R₅ and R₆ in each case denote a hydrogen atom, and

R₇ is a ~~fluorine, chlorine, bromine or iodine atom;~~

a methoxy, nitro, cyano, carboxy, C₁₋₃-alkoxycarbonyl, aminocarbonyl, C₁₋₃-alkylaminocarbonyl, di-(C₁₋₃-alkyl) aminocarbonyl, phenyl C₁₋₃-alkylaminocarbonyl, N-(phenyl C₁₋₃-alkyl) C₁₋₃-alkylaminocarbonyl or 5- to 7-membered cycloalkyleneiminocarbonyl group,
a C₁₋₃-alkyl group which is substituted by a carboxy, C₁₋₃-alkoxycarbonyl, aminocarbonyl, C₁₋₃-alkylaminocarbonyl, di-(C₁₋₃-alkyl) aminocarbonyl, phenyl C₁₋₃-alkylaminocarbonyl, N-(phenyl C₁₋₃-alkyl) C₁₋₃-alkylaminocarbonyl, 5- to 7-membered cycloalkyleneiminocarbonyl, amino, C₁₋₃-alkylamino, di-(C₁₋₃-alkyl) amino, phenyl C₁₋₃-alkylamino, N-(phenyl C₁₋₃-alkyl) C₁₋₃-alkylamino or 5- to 7-membered cycloalkyleneimino piperidino group,
while the abovementioned 5- to 7-membered cycloalkyleneimino piperidino group may be substituted by one or two C₁₋₃-alkyl groups, which may in turn be terminally substituted by a hydroxy, amino or C₂₋₄-alkanoylamino group, and at the same time in the abovementioned piperidino group 5- to 7-membered cycloalkyleneimino moieties a methylene group in the 2 position may be replaced by a carbonyl group or in the abovementioned 6- and 7-membered cycloalkyleneimino moieties a methylene group in the 4 position may be replaced by an oxygen atom, by an imino, N-(C₁₋₃-alkyl) imino, N-(phenyl C₁₋₃-alkyl) imino or N-(C₁₋₅-alkoxycarbonyl) imino group,
an amino, C₁₋₃-alkylamino, phenyl C₁₋₃-alkylamino, C₁₋₅-alkanoylamino, phenyl C₁₋₄-alkanoylamino, C₁₋₅-alkoxycarbonylamino, phenyl C₁₋₃-alkoxycarbonylamino, C₁₋₅-alkylsulphonylamino, phenyl C₁₋₃-alkylsulphonylamino or phenylsulphonylamino group,
wherein the hydrogen atom of the amino group may be replaced by a C₁₋₃-alkyl group, while the C₁₋₃-alkyl moiety may be substituted by a carboxy, C₁₋₃-alkoxycarbonyl, aminocarbonyl, C₁₋₃-alkylaminocarbonyl, di-(C₁₋₃-alkyl) aminocarbonyl, phenyl C₁₋₃-alkylaminocarbonyl, N-(phenyl C₁₋₃-alkyl) C₁₋₃-alkylaminocarbonyl, 2-dimethylaminoethylaminocarbonyl, N-methyl (2-dimethylaminoethyl) aminocarbonyl or C₄₋₆-cycloalkyleneiminocarbonyl group or from position 2 by an amino, C₁₋₃-alkylamino, di-(C₁₋₃-alkyl) amino, phenyl C₁₋₃-alkylamino, N-(phenyl C₁₋₃-alkyl) C₁₋₃-alkylamino, C₂₋₅-alkanoylamino, N-(C₁₋₃-alkyl) C₂₋₅-alkanoylamino, C₁₋₅-alkoxycarbonylamino or N-(C₁₋₅-alkoxycarbonyl) C₁₋₃-alkylamino group.

Claim 6 (currently amended): A compound of formula I according to claim 1, wherein:

R₂ is a C₁₋₃-alkyl group optionally substituted by a phenyl group, a C₁₋₃-perfluoroalkyl group or a phenylvinyl group, or
a phenyl group which may be substituted by a fluorine, chlorine, bromine or iodine atom, by a C₁₋₃-alkyl, C₁₋₃-alkoxy, nitro, amino, cyano, ~~cyanomethyl~~ or aminomethyl group, a C₄₋₆-alkyl, C₃₋₇-cycloalkyl, trimethylphenyl or naphthyl group,
~~a pyridinyl, quinolyl, isoquinolyl, oxazolyl, isoxazolyl, imidazolyl or 1-(C₁₋₃-alkyl)-imidazolyl group optionally substituted by a C₁₋₃-alkyl group,~~

R₃ is a phenyl group optionally substituted by a fluorine, chlorine, bromine or iodine atom, by a C₁₋₃-alkyl, C₁₋₃-alkoxy, nitro or amino group,

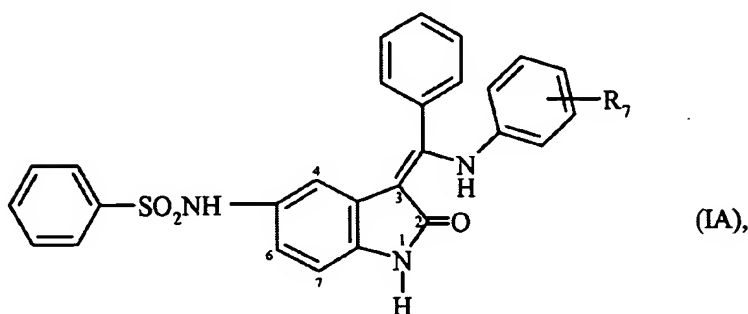
R₄ is a phenyl group which ~~may be~~ is substituted by R₇ and additionally by a chlorine atom or a nitro group, while

R₇ is ~~a fluorine, chlorine, bromine or iodine atom,~~
~~a methoxy, nitro, cyano, carboxy, methoxycarbonyl, aminocarbonyl, methylaminocarbonyl, dimethylaminocarbonyl, benzylaminocarbonyl, N-benzyl-methylaminocarbonyl, pyrrolidinocarbonyl or piperidinocarbonyl group,~~
a methyl or ethyl group which ~~may be~~ is substituted by a ~~carboxy, methoxycarbonyl, aminocarbonyl, methylaminocarbonyl, dimethylaminocarbonyl, benzylaminocarbonyl, N-benzyl-methylaminocarbonyl, pyrrolidinocarbonyl, piperidinocarbonyl, amino, methylamino, dimethylamino, benzylamino, N-benzylmethylamino, C₂₋₄-alkanoylamino, N-methyl-C₂₋₄-alkanoylamino, tert.butyloxycarbonylamino, N-methyl-tert.butyloxycarbonylamino, pyrrolidino, pyrrolidinomethyl, hydroxypyrrolidinomethyl, hydroxymethylpyrrolidinomethyl, piperidino, dimethylpiperidino, 2-oxo-piperidino, piperazino, 4-methyl-piperazino, 4-benzyl-piperazino, 4-tert.butoxycarbonyl-piperazino or morpholine group, or~~
~~an amino, methylamino, ethylamino, C₁₋₃-alkanoylamino, phenylacetylamino, tert.butoxycarbonylamino, C₁₋₄-alkylsulphonylamino, phenyl-methylsulphonylamino or phenylsulphonylamino group, wherein the hydrogen atom of the amino group may be replaced by a methyl or ethyl group, while the methyl or ethyl moiety in each case may be substituted by a carboxy, methoxycarbonyl, aminocarbonyl, methylaminocarbonyl or dimethylaminocarbonyl group or the ethyl moiety may also be substituted from position 2 by~~

an amino, methylamino, dimethylamino, benzylalkylamino, N benzyl methylamino, C₂₋₃-alkanoylamino, N-methyl-C₂₋₃-alkanoylamino, tert.butyloxycarbonylamino or N-methyl-tert.butyloxycarbonylamino group.

Claim 7 (original): A compound of formula I according to claim 1, wherein R₄ is a phenyl group substituted in the 4 position by R₇.

Claim 8 (original): A compound of the formula IA



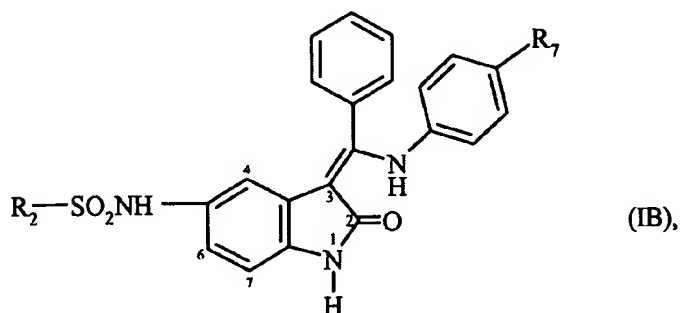
wherein R₇ is defined as in claim 1, 5 or 6.

Claim 9 (original): A compound of formula IA according to claim 8 wherein R₇ is selected from the group consisting of:

hydrogen, (2,6-dimethylpiperidino)-methyl, (N-ethylsulphonyl)-N-(2-dimethylaminoethyl)-aminocarbonylmethyl-amino, N-ethylsulphonyl-N-(N-(2-dimethylaminoethyl)-N-methyl-amino-carbonylmethyl)-amino, and 2-oxopiperidinomethyl, 4-benzyl-piperazino-methyl, 4-methylpiperazino-methyl, 4-tert.butyloxycarbonyl-piperazinomethyl, acetylamino, acetylaminomethyl, amino, aminomethyl, benzylaminocarbonyl, benzylaminocarbonylmethyl, carboxy, carboxymethyl, chlorine, cyano, dimethylaminocarbonyl-methylamino, dimethylaminoethyl, dimethylaminomethyl, ethoxy-carbonylmethyl, ethylsulphonylamino, formylamino, methoxy-carbonyl, methylsulphonylamino, morpholinomethyl, N-(2-(N-acetyl-N-methyl-amino)-ethyl)-ethylsulphonylamino, N-(2-(N-acetyl-N-methyl-amino)-ethyl)-methylsulphonylamino, N-(2-(N-acetyl-N-methyl-amino)-ethyl)-propionylamino, N-(2-(N-acetyl-N-methyl-amino)-ethyl)amino, N-(2-(N-benzyl-N-methyl-amino)-ethyl)-propionyl-amino, N-(2-acetyl-amino-ethyl)-N-acetyl-amino, N-(2-acetyl-amino-ethyl)-N-ethylsulphonyl-

~~amino, N (2 acetyl amino ethyl) N methylsulphonyl amino, N (2 acetyl amino ethyl) N propionyl amino, N (2 aminoethyl) N methylsulphonyl amino, N (2 dimethylamino ethyl) N acetyl amino, N (2 dimethylamino ethyl) N butylsulphonyl amino, N (2 dimethylamino ethyl) N methylsulphonyl amino, N (2 dimethylamino ethyl) N phenylsulphonyl amino, N (2 dimethylaminoethyl) N propylsulphonyl amino, N (2 methylamino ethyl) acetyl amino, N (2 methylamino ethyl) N methylsulphonyl amino, N (2 methylamino ethyl) propionyl amino, N (2 propionyl amino ethyl) N propionyl amino, N (aminocarbonyl methyl) N methylsulphonyl amino, N (dimethylamino carbonylmethyl) N (methylsulphonyl amino, N (dimethylaminoethyl) N methylsulphonyl amino, N (methylaminocarbonyl methyl) N methylsulphonyl amino, N (piperidinomethyl carbonyl) N methyl amino, N acetyl N (2 (N benzyl N methyl amino) ethyl amino, N acetyl N (2 benzyl oxycarbonylamino ethyl) amino, N carboxylmethyl N methylsulphonyl amino, N ethylsulphonyl N hydroxycarbonylmethyl amino, N methyl N acetyl amino, N methyl N ethylsulphonyl amino, N methyl N formyl amino, N methyl N methylsulphonyl amino, N methyl N propionyl amino, piperazinomethyl, propionyl amino, pyrrolidin 1 yl methyl, 2 hydroxymethylpyrrolidin 1 yl methyl, 3 hydroxypyrrolidin 1 yl methyl and tert.butoxycarbonylamino.~~

Claim 10 (original): A compound of formula IB



wherein R₂ and R₇ are defined as in claim 1, 4, 5 or 6.

Claim 11 (currently amended): A compound of formula IB according to claim 10 wherein:

R₇ is selected from the group consisting of:

hydrogen, (2,6-dimethylpiperidino)-methyl ~~and, (N-ethylsulphonyl)-N-(2-~~
~~dimethylaminoethyl)-aminocarbonylmethyl)-amino, N-ethylsulphonyl-N-(2-~~
~~dimethylaminoethyl)-N-methyl-amino-carbonylmethyl)-amino, 2-oxopiperidinomethyl, 4-~~
~~benzyl-piperazino-methyl, 4-methylpiperazino-methyl, 4-tert.butoxycarbonyl-~~
~~piperazinomethyl, acetylamino, acetylaminomethyl, amino, aminomethyl,~~
~~benzylaminocarbonyl, benzylaminocarbonyl-methyl, carboxy, carboxymethyl, chlorine,~~
~~cyano, dimethylaminocarbonyl-methylamino, dimethylaminoethyl, dimethylaminomethyl,~~
~~ethoxycarbonylmethyl, ethylsulphonylamino, formylamino, methoxycarbonyl,~~
~~methylsulphonylamino, morpholinomethyl, N-(2-(N-acetyl-N-methyl-amino)-ethyl)-~~
~~ethylsulphonylamino, N-(2-(N-acetyl-N-methyl-amino)-ethyl)-methylsulphonylamino, N-(2-~~
~~(N-acetyl-N-methyl-amino)-ethyl)-propionylamino, N-(2-(N-acetyl-N-methyl-amino)-~~
~~ethylamino, N-(2-(N-benzyl-N-methyl-amino)-ethyl)-propionylamino, N-(2-acetyl-amino-~~
~~ethyl)-N-acetyl-amino, N-(2-acetyl-amino-ethyl)-N-ethylsulphonyl-amino, N-(2-acetyl-amino-~~
~~ethyl)-N-methylsulphonyl-amino, N-(2-acetyl-amino-ethyl)-N-propionyl-amino, N-(2-~~
~~aminoethyl)-N-methylsulphonyl-amino, N-(2-dimethyl-amino-ethyl)-N-acetyl-amino, N-(2-~~
~~dimethyl-amino-ethyl)-N-butylsulphonyl-amino, N-(2-dimethyl-amino-ethyl)-N-~~
~~methylsulphonyl-amino, N-(2-dimethyl-amino-ethyl)-N-phenylsulphonyl-amino, N-(2-~~
~~dimethylaminoethyl)-N-propylsulphonyl-amino, N-(2-methyl-amino-ethyl)-acetyl-amino, N-~~
~~(2-methyl-amino-ethyl)-N-methylsulphonyl-amino, N-(2-methyl-amino-ethyl)-~~
~~propionylamino, N-(2-propionyl-amino-ethyl)-N-propionyl-amino, N-(aminocarbonyl-~~
~~methyl)-N-methylsulphonyl-amino, N-(dimethyl-amino-carbonylmethyl)-N-~~
~~(methylsulphonyl-amino, N-(dimethylaminoethyl)-N-methylsulphonyl-amino, N-~~
~~(methylaminocarbonyl-methyl)-N-methylsulphonyl-amino, N-(piperidinomethyl-carbonyl)-~~
~~N-methyl-amino, N-acetyl-N-(2-(N-benzyl-N-methyl-amino)-ethylamino, N-acetyl-N-(2-~~
~~benzyl-oxycarbonylamino-ethyl)-amino, N-carboxylmethyl-N-methylsulphonyl-amino, N-~~
~~ethylsulphonyl-N-hydroxycarbonylmethyl-amino, N-methyl-N-acetyl-amino, N-methyl-N-~~
~~ethylsulphonyl-amino, N-methyl-N-formyl-amino, N-methyl-N-methylsulphonyl-amino, N-~~
~~methyl-N-propionyl-amino, piperazinomethyl, propionylamino, pyrrolidin-1-yl-methyl, 2-~~

~~hydroxymethylpyrrolidin-1-yl-methyl, 3-hydroxypyrrolidin-1-yl-methyl and
tert.butoxycarbonylamino; and~~

R₂ is selected from the group consisting of:

~~1-methyl-1H-imidazol-4-yl, 2-aminophenyl, 2-chlorophenyl, 2-cyanophenyl, 2-nitrophenyl,
2-phenylethene, 3-aminomethylphenyl, 3-aminophenyl, 3-chlorophenyl, 3-cyanophenyl, 3-
methoxyphenyl, 3-methylphenyl, 3-nitrophenyl, 4-aminophenyl, 4-chlorophenyl,
4-methoxyphenyl, 4-methylphenyl, 4-nitrophenyl, benzyl, quinolin-8-yl, cyclopropyl, ethyl,
isopropyl, methyl, naphthalin-1-yl, naphthalin-2-yl, propyl, pyrid-2-yl, pyrid-3-yl, 3,5-
dimethyl-isoxazol-4-yl and 2,4,6-trimethylphenyl.~~

Claim 12 (currently amended): A compound selected from the group consisting of:

~~(Z)-3-{1-[4-(N-(2-aminoethyl)-N-methylsulphonylamino)-phenylamino]-1-phenyl-
methylidene}-5-phenylsulphonylamino-2-indolinone;
(Z)-3-{1-[4-(N-(2-dimethylaminoethyl)-N-phenylsulphonylamino)-phenylamino]-1-phenyl-
methylidene}-5-phenylsulphonylamino-2-indolinone;
(Z)-3-{1-[4-(4-methylpiperazinomethyl)-phenylamino]-1-phenyl-methylidene}-5-
phenylsulphonylamino-2-indolinone;
(Z)-3-{1-[4-(pyrrolidin-1-ylmethyl)-phenylamino]-1-phenyl-methylidene}-5-
phenylsulphonylamino-2-indolinone;
(Z)-3-{1-[4-(N-methyl-N-acetyl-amino)-phenylamino]-1-phenyl-methylidene}-5-
phenylsulphonylamino-2-indolinone;
(Z)-3-(1-phenylamino-1-phenyl-methylidene)-5-phenylsulphonylamino-2-indolinone;
(Z)-3-[1-(4-chlorophenylamino)-1-phenyl-methylidene]-5-phenylsulphonylamino-2-
indolinone;
(Z)-3-{1-[4-(N-(2-propionylamino-ethyl)-N-propionyl-amino)-phenylamino]-1-phenyl-
methylidene}-5-phenylsulphonylamino-2-indolinone;
(Z)-3-[1-(4-dimethylaminomethyl-phenylamino)-1-phenyl-methylidene]-5-
phenylsulphonylamino-2-indole;~~

~~(Z)-3-[1-(4-(N-methyl-N-methylsulphonyl-amino)-phenylamino)-1-phenyl-methylidene]-5-phenylsulphonylamino-2-indolinone,~~
~~(Z)-3-[1-(4-(N-methyl-N-piperidinomethylcarbonyl-amino)-phenylamino)-1-phenyl-methylidene]-5-phenylsulphonylamino-2-indolinone,~~
~~(Z)-3-[1-(4-(pyrrolidin-1-ylmethyl)-phenylamino)-1-phenyl-methylidene]-5-benzylsulphonylamino-2-indolinone,~~
~~(Z)-3-[1-(4-((2,6-dimethylpiperidino)-methyl)-phenylamino)-1-phenyl-methylidene]-5-(3-nitrophenylsulphonylamino)-2-indolinone,~~
~~(Z)-3-[1-(4-dimethylaminomethyl-phenylamino)-1-phenyl-methylidene]-5-ethylsulphonylamino-2-indolinone,~~
~~(Z)-3-[1-(4-(N-benzyl-N-methyl-aminomethyl)-phenylamino)-1-phenyl-methylidene]-5-ethylsulphonylamino-2-indolinone,~~
~~(Z)-3-[1-(4-(2-dimethylamino-ethyl)-phenylamino)-1-phenyl-methylidene]-5-ethylsulphonylamino-2-indolinone,~~
~~(Z)-3-[1-(4-(pyrrolidin-1-ylmethyl)-phenylamino)-1-phenyl-methylidene]-5-(pyridin-3-ylsulphonylamino)-2-indolinone,~~
~~(Z)-3-[1-(4-(pyrrolidin-1-ylcarbonyl)-phenylamino)-1-phenyl-methylidene]-5-(pyridin-3-ylsulphonylamino)-2-indolinone,~~
~~(Z)-3-[1-(4-(piperidinomethyl)-phenylamino)-1-phenyl-methylidene]-5-methylsulphonylamino-2-indolinone,~~
~~(Z)-3-[1-(4-(piperidinomethyl)-phenylamino)-1-phenyl-methylidene]-5-ethylsulphonylamino-2-indolinone,~~
~~(Z)-3-[1-(4-(piperidinomethyl)-phenylamino)-1-phenyl-methylidene]-5-isopropylsulphonylamino-2-indolinone,~~
~~(Z)-3-[1-(4-(piperidinomethyl)-phenylamino)-1-phenyl-methylidene]-5-(naphthalin-1-ylsulphonylamino)-2-indolinone,~~
~~(Z)-3-[1-(4-(piperidinomethyl)-phenylamino)-1-phenyl-methylidene]-5-(3-nitrophenylsulphonylamino)-2-indolinone,~~ and
~~(Z)-3-[1-(4-(piperidinomethyl)-phenylamino)-1-phenyl-methylidene]-5-(3,5-dimethylisoxazol-4-ylsulphonylamino)-2-indolinone,~~

~~(Z)-3-[1-[4-(piperidinomethyl)-phenylamino]-1-phenyl-methylidene]-5-cyclopropylsulphonylamino-2-indolinone;~~
~~(Z)-3-[1-[4-(piperidinomethyl)-phenylamino]-1-phenyl-methylidene]-5-(pyridin-3-ylphenylsulphonylamino)-2-indolinone;~~
~~(Z)-3-[1-[4-(pyrrolidin-1-ylmethyl)-phenylamino]-1-phenyl-methylidene]-5-cyclopropylsulphonylamino-2-indolinone;~~
~~(Z)-3-[1-[4-(pyrrolidin-1-ylmethyl)-phenylamino]-1-phenyl-methylidene]-5-propylsulphonylamino-2-indolinone;~~
~~(Z)-3-[1-[4-(pyrrolidin-1-ylmethyl)-phenylamino]-1-phenyl-methylidene]-5-ethylsulphonylamino-2-indolinone;~~
~~(Z)-3-[1-[4-(pyrrolidin-1-ylmethyl)-phenylamino]-1-phenyl-methylidene]-5-methylsulphonylamino-2-indolinone;~~
~~(Z)-3-[1-[4-(benzylaminocarbonyl)-phenylamino]-1-phenyl-methylidene]-5-phenylsulphonylamino-2-indolinone;~~
~~(Z)-3-[1-[4-(N-dimethylaminocarbonylmethyl-N-acetyl-amino)-phenylamino]-1-phenyl-methylidene]-5-phenylsulphonylamino-2-indolinone;~~
(Z)-3-[1-(4-piperidinomethyl-phenylamino)-1-phenyl-methylidene]-5-(4-aminophenylsulphonylamino)-2-indolinone, and
~~(Z)-3-[1-[4-(N-(2-dimethylamino-ethyl)-N-methylsulphonyl-amino)-phenylamino]-1-phenyl-methylidene]-5-(N-methyl-N-phenylsulphonyl-amino)-2-indolinone;~~
or a pharmaceutically acceptable salt thereof.

Claim 13 (currently amended): A pharmaceutical preparation comprising a compound according to claim 1, 2, 3, 4, 5, 6, 7, ~~8, 9, 10, 11~~ or 12 and a pharmaceutically acceptable carrier.

Claim 14 (withdrawn): A method for treating a disease characterised by excessive or abnormal cell proliferation which comprises administering a therapeutic amount of a compound according to claim 1, 2, 3, 4, 5, 6, 7, ~~8, 9, 10, 11~~ or 12.